

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121318\
 Data File : BE098414.D
 Acq On : 14 Dec 2018 1:44
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 14 02:15:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 14:59:55 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	84	0.00
2	1,4-Dioxane	0.692	0.641	7.4	79	-0.01
3	n-Nitrosodimethylamine	0.623	0.369	40.8#	54	0.00
4 S	2-Fluorophenol	0.936	0.883	5.7	82	0.00
5 S	Phenol-d6	1.326	1.271	4.1	80	0.00
6	bis(2-Chloroethyl)ether	1.260	1.089	13.6	75	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	83	-0.01
8 S	Nitrobenzene-d5	0.364	0.347	4.7	86	0.01
9	Naphthalene	4.025	3.837	4.7	81	0.00
10	Hexachlorobutadiene	0.256	0.272	-6.3	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.650	0.591	9.1	76	0.00
12	2-Methylnaphthalene	0.654	0.584	10.7	76	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
14 S	2,4,6-Tribromophenol	0.147	0.116	21.1	69	0.00
15 S	2-Fluorobiphenyl	1.687	1.496	11.3	80	0.00
16	Acenaphthylene	1.874	1.790	4.5	81	0.00
17	Acenaphthene	1.126	1.024	9.1	75	0.00
18	Fluorene	1.506	1.397	7.2	77	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	79	-0.01
20	4-Bromophenyl-phenylether	0.215	0.186	13.5	71	0.00
21	Hexachlorobenzene	0.231	0.217	6.1	75	0.00
22	Pentachlorophenol	0.046	0.014	69.6#	30#	0.01
23	Phenanthrene	0.972	0.908	6.6	75	0.00
24	Anthracene	0.866	0.735	15.1	68	0.00
25 SURR	Fluoranthene-d10	5.131	4.841	5.7	72	0.00
26	Fluoranthene	1.286	1.249	2.9	74	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	88	-0.01
28	Pvrene	1.255	1.088	13.3	75	0.00
29 S	Terphenyl-d14	0.972	0.773	20.5	71	0.00
30	Benzo(a)anthracene	1.139	1.042	8.5	82	-0.01
31	Chrysene	1.193	1.111	6.9	83	0.00
32	Bis(2-ethylhexyl)phthalate	2.313	1.680	27.4#	64	-0.01
33	Indeno(1,2,3-cd)pyrene	1.187	1.173	1.2	92	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	93	-0.01
35	Benzo(b)fluoranthene	1.094	1.076	1.6	95	-0.01
36	Benzo(k)fluoranthene	1.274	1.342	-5.3	102	-0.01
37 C	Benzo(a)pyrene	1.129	1.113	1.4	94	-0.01
38	Dibenzo(a,h)anthracene	1.063	0.989	7.0	89	-0.01
39	Benzo(a,h,i)perylene	1.071	1.044	2.5	94	-0.02

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0			